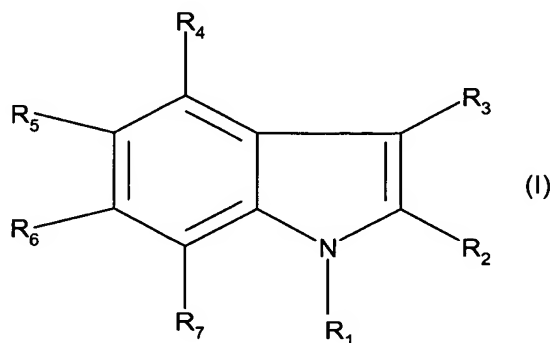


### Amendments to the Claims

This listing of claims will replace all prior versions, and listings, of claims in the application:

#### Listing of Claims:

1. (Previously Presented) An indole compound represented by the formula (I), or a pharmaceutically acceptable salt, solvate, or prodrug thereof;



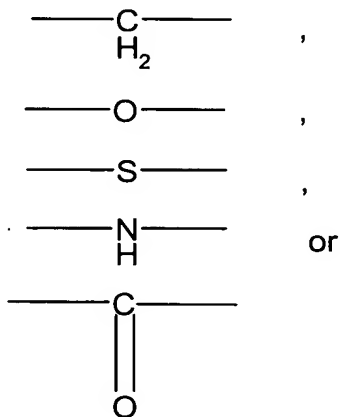
wherein ;

$R_1$  is (c) wherein;

(c) is the group  $-(L_1)-R_{11}$ ; where,  $-(L_1)-$  is a divalent linking group of 1 to 8 atoms and where  $R_{11}$  is  $-(CH_2)_m-R_{12}$ ;

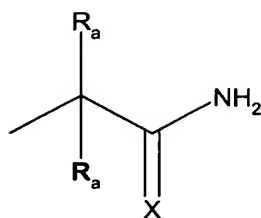
$R_2$  is hydrogen, or C1-C4 alkyl;

$R_3$  is  $-(L_3)-Z$ , where  $-(L_3)-$  is a divalent linker group selected from a bond or:



$-\text{CH}_2-$

and Z is a group represented by the formulae,



wherein, X is oxygen or sulfur; and R<sub>a</sub> is selected from hydrogen, C<sub>1</sub>-C<sub>8</sub> alkyl, aryl, C<sub>1</sub>-C<sub>8</sub> alkaryl, C<sub>1</sub>-C<sub>8</sub> alkoxy, aralkyl and -CN;

R<sub>4</sub> is the group, -(L<sub>h</sub>)-(hydroxyfunctional amide); wherein -(L<sub>h</sub>)-, is an hydroxyfunctional amide linker having an hydroxyfunctional amide linker length of 1 to 8;

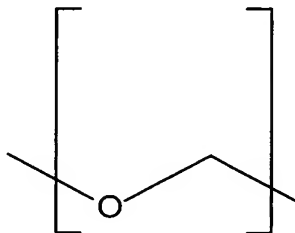
R<sub>5</sub> is selected from hydrogen, a non-interfering substituent, or the group, -(L<sub>a</sub>)-(acidic group); wherein -(L<sub>a</sub>)-, is an acid linker having an acid linker length of 1 to 8;

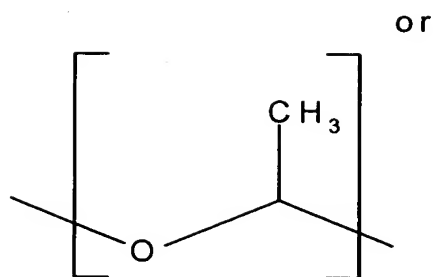
R<sub>6</sub> and R<sub>7</sub> are selected from hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>2</sub>-C<sub>6</sub> alkenyl, and C<sub>2</sub>-C<sub>6</sub> alkynyl.

2. (Previously Presented) The compound of claim 1 wherein R<sub>2</sub> is hydrogen, C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>2</sub>-C<sub>4</sub> alkenyl, -O-(C<sub>1</sub>-C<sub>3</sub> alkyl), -S-(C<sub>1</sub>-C<sub>3</sub> alkyl), and C<sub>3</sub>-C<sub>4</sub> cycloalkyl.

3. (Cancelled)

4. (Previously Presented) The compound of Claim 1 wherein the hydroxyfunctional amide linker group, -(L<sub>h</sub>)-, for R<sub>4</sub> is a divalent group selected from,





where R<sub>40</sub>, R<sub>41</sub>, R<sub>42</sub>, and R<sub>43</sub> are each independently selected from hydrogen, C<sub>1</sub>-C<sub>8</sub> alkyl.

5. (Cancelled)

6. (Previously Presented) The compound of claim 1 wherein R<sub>5</sub> is the group, -(L<sub>a</sub>)-(acidic group) and wherein the {acidic group} is:

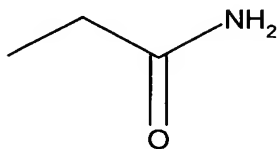


7. (Cancelled)

8. (Cancelled)

9. (Cancelled)

10. (Original) The compound of claim 1 wherein for R<sub>3</sub>, Z is the group represented by the formula;



and the linking group -(L<sub>3</sub>)- is a bond.

11. (Cancelled)

12. (Cancelled)

13. (Cancelled)

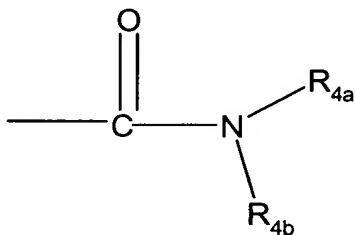
14. (Cancelled)

15. (Cancelled)

16. (Cancelled)

17. (Cancelled)

18. (Original) The compound of claim 1 wherein R<sub>4</sub> is the group, -(L<sub>C</sub>)-(hydroxyfunctional amide group) and wherein the (hydroxyfunctional amide group) is:



and R<sup>4a</sup> is independently selected from the group consisting of OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, (C<sub>7</sub>-C<sub>14</sub>)alkaryloxy, (C<sub>2</sub>-C<sub>8</sub>)alkenyloxy, (C<sub>7</sub>-C<sub>14</sub>) aralkyloxy, (C<sub>7</sub>-C<sub>14</sub>)aralkenyloxy and aryloxy; and

wherein R<sup>4b</sup> is independently selected from the group consisting of H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, arylalkyl, heteroaryl and aryl.

19. (Cancelled)

20. (Previously Presented) A compound selected from the group of:

2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(hydroxy)acetamide;

2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(methoxy)acetamide;

2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(methyl)-N-(methoxy)acetamide;

2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(hydroxy)-N-(methyl)acetamide;

2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(ethoxy)acetamide;

2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(2-propenyloxy)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(hydroxy)-*N*-(2-propyl)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(tert-butyloxy)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-[2-(methyl)propyloxy]acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(phenylmethyloxy)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(methyl)-*N*-(phenylmethyloxy)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(phenyloxy)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(methyl)-*N*-(phenyloxy)acetamide;  
2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(cyclohexyl)-*N*-(hydroxy)acetamide; and  
2-[[3-(2-Amino-2-oxoethyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(hydroxy)acetamide.

21. (Cancelled)

22. (Original) A pharmaceutical formulation comprising a indole compound as claimed in claim 1 together with a pharmaceutically acceptable carrier or diluent therefor.

23. (Cancelled)

24. (Cancelled)

25. (Previously Presented) A pharmaceutical formulation containing an effective amount of the compound of claim 1 useful for the treatment and/or amelioration of Inflammatory Diseases.

26. (Cancelled)

Serial No. 10/629,992

27. (Cancelled)